

Refine Search

Search Results -

Term	Documents
(7 AND 18).PGPB,USPT,USOC,EPAB,JPAB,DWPI,TDBD.	0
(L18 AND L7).PGPB,USPT,USOC,EPAB,JPAB,DWPI,TDBD.	0

Database:

US Pre-Grant Publication Full-Text Database
 US Patents Full-Text Database
 US OCR Full-Text Database
 EPO Abstracts Database
 JPO Abstracts Database
 Derwent World Patents Index
 IBM Technical Disclosure Bulletins

Search:

L20

Refine Search

Recall Text

Clear

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Search History

DATE: Monday, August 15, 2005 [Printable Copy](#) [Create Case](#)

Set Name **Query**
 side by side

Hit Count **Set Name**
 result set

DB=PGPB,USPT,USOC,EPAB,JPAB,DWPI,TDBD; PLUR=YES; OP=ADJ

<u>L20</u>	118 and 17	0	<u>L20</u>
<u>L19</u>	L18 and 18	1	<u>L19</u>
<u>L18</u>	indigo carmine	526	<u>L18</u>
<u>L17</u>	L16 and 18	2	<u>L17</u>
<u>L16</u>	indigo	10429	<u>L16</u>
<u>L15</u>	skin same 18	7	<u>L15</u>
<u>L14</u>	skin and 18	648	<u>L14</u>
<u>L13</u>	hair and 18	32	<u>L13</u>
<u>L12</u>	110 same 18	1	<u>L12</u>
<u>L11</u>	110 and 18	541	<u>L11</u>
<u>L10</u>	cosmetic	171283	<u>L10</u>
<u>L9</u>	18 and 17	2	<u>L9</u>
<u>L8</u>	dehydrocholic acid	1022	<u>L8</u>

<u>L7</u>	indigofera tinctoria	29	<u>L7</u>
<u>L6</u>	L5 and l2	8	<u>L6</u>
<u>L5</u>	l4 or l1	61	<u>L5</u>
<u>L4</u>	patcholi	2	<u>L4</u>
<u>L3</u>	l1 and l2	8	<u>L3</u>
<u>L2</u>	n-hexane or chloroform or n-butanol	246065	<u>L2</u>
<u>L1</u>	agastache rugosa	60	<u>L1</u>

END OF SEARCH HISTORY

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NEWS	4	FEB 28	BABS - Current-awareness alerts (SDIs) available
NEWS	5	MAR 02	GBFULL: New full-text patent database on STN
NEWS	6	MAR 03	REGISTRY/ZREGISTRY - Sequence annotations enhanced
NEWS	7	MAR 03	MEDLINE file segment of TOXCENTER reloaded
NEWS	8	MAR 22	KOREAPAT now updated monthly; patent information enhanced
NEWS	9	MAR 22	Original IDE display format returns to REGISTRY/ZREGISTRY
NEWS	10	MAR 22	PATDPASPC - New patent database available
NEWS	11	MAR 22	REGISTRY/ZREGISTRY enhanced with experimental property tags
NEWS	12	APR 04	EPPFULL enhanced with additional patent information and new fields
NEWS	13	APR 04	EMBASE - Database reloaded and enhanced
NEWS	14	APR 18	New CAS Information Use Policies available online
NEWS	15	APR 25	Patent searching, including current-awareness alerts (SDIs), based on application date in CA/CAPLUS and USPATFULL/USPAT2 may be affected by a change in filing date for U.S. applications.
NEWS	16	APR 28	Improved searching of U.S. Patent Classifications for U.S. patent records in CA/CAPLUS
NEWS	17	MAY 23	GBFULL enhanced with patent drawing images
NEWS	18	MAY 23	REGISTRY has been enhanced with source information from CHEMCATS
NEWS	19	JUN 06	The Analysis Edition of STN Express with Discover! (Version 8.0 for Windows) now available
NEWS	20	JUN 13	RUSSIAPAT: New full-text patent database on STN
NEWS	21	JUN 13	FRFULL enhanced with patent drawing images
NEWS	22	JUN 27	MARPAT displays enhanced with expanded G-group definitions and text labels
NEWS	23	JUL 01	MEDICONF removed from STN
NEWS	24	JUL 07	STN Patent Forums to be held in July 2005
NEWS	25	JUL 13	SCISEARCH reloaded
NEWS	26	JUL 20	Powerful new interactive analysis and visualization software, STN AnaVist, now available
NEWS	27	AUG 11	Derwent World Patents Index(R) web-based training during August
NEWS	28	AUG 11	STN AnaVist workshops to be held in North America
NEWS EXPRESS			JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005
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NEWS INTER			General Internet Information
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NEWS PHONE			Direct Dial and Telecommunication Network Access to STN
NEWS WWW			CAS World Wide Web Site (general information)

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 15:46:04 ON 15 AUG 2005

=> file ca, biosis, medline

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FILE 'BIOSIS' ENTERED AT 15:46:17 ON 15 AUG 2005

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FILE 'MEDLINE' ENTERED AT 15:46:17 ON 15 AUG 2005

=> s agastache rugosa or patcholi

L1 161 AGASTACHE RUGOSA OR PATCHOLI

=> s n-hexane or chloroform or n-butanol

L2 108333 N-HEXANE OR CHLOROFORM OR N-BUTANOL

=> s l1 and l2

L3 3 L1 AND L2

=> d 1-3 ab,bib

L3 ANSWER 1 OF 3 CA COPYRIGHT 2005 ACS on STN

AB A novel compound and its pharmaceutically acceptable salt separated from **Agastache rugosa**, its preparation method and a pharmaceutical composition containing the compound and/or the salt are provided, which compound is used as an apoptosis inhibitor. The novel compound is represented by the formula 1 or 2. In the formulas 1 and 2, R1, R2, R3, R4 and R5 are independent each another and are an alkylalkoxy group of C1-C4; H; a straight or branched alkyl group of C1-C6; OH; a dialkylamino group of C2-C6; a straight or branched hydroxyalkyl group of C1-C6; a straight or branched dihydroxyalkyl group of C3-C6; an alkoxyalkyl group of C3-C6; or a saturated or unsatd. five or six-membered hetero ring containing 1-3 hetero atoms selected from N, O and S and unsubstituted or substituted with an alkyl group of C1-C3. The method comprises the steps of extracting **Agastache rugosa** with an organic solvent 1, concentrating the extract under reduced pressure and extracting the concentrated one with an organic solvent 2 to obtain the extract; concentrating the extract, dissolving the concentrated one into a small quantity of a solvent mixture, and obtaining the concentrated active fraction by chromatog.; dissolving the concentrated fraction into an organic solvent 3 and obtaining an active fraction by chromatog.; and separating the active fraction by chromatog. of the obtained fraction, removing the solvent and lyophilizing it. Preferably the organic solvents 1 and 3 are selected from the group consisting of methanol, ethanol, acetone and distilled water; the organic solvent 2 is selected from the group consisting of Et acetate, hexane and butanol; and the solvent mixture is a mixture of **chloroform** and methanol or hexane and Et acetate.

AN 142:110549 CA

TI Method for preparation of novel compound separated from **Agastache rugosa** with apoptosis inhibiting activity as pharmaceutical composition

IN Kim, Hyeon A.; Ko, Yeong Hui; Lee, Chung Hwan; Lee, Ho Jae; Lee, Sang Myeong

PA Korea Research Institute of Bioscience and Biotechnology, S. Korea
SO Repub. Korean Kongkae Taeho Kongbo, No pp. given
CODEN: KRXXA7
DT Patent
LA Korean
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	KR 2002059986	A	20020716	KR 2001-1180	20010109
PRAI	KR 2001-1180		20010109		

L3 ANSWER 2 OF 3 CA COPYRIGHT 2005 ACS on STN

AB Thirty plants were screened for their antioxidative activity. The exts. of **Agastache rugosa**, *Schizonepeta tenuifolia* and *Lycopus lucidus* had high free radical (2,2-diphenyl-1-picrylhydrazyl) scavenging activities. Methanol extract of **Agastache rugosa** was fractionated with hexane, **chloroform**, Et acetate, and water. The Et acetate fraction showed the highest antioxidant activity in the DPPH test. The Et acetate fraction was applied to Sephadex LH-20 column, and the fractions showing antioxidative activity were collected and used for identification of the substance. The purified substance was applied to mass, IR, UV and NMR spectroscopy. The spectra of mass, IR, UV and NMR implied that the substance was a phenolic compound rosmarinic acid. The rosmarinic acid had more antioxidative effect than those of BHT and α -tocopherol in the Rancimat test.

AN 132:205431 CA

TI Isolation, identification, and activity of rosmarinic acid, a potent antioxidant extracted from Korean **Agastache rugosa**

AU Kim, Jung-Bong; Kim, Jong-Bum; Cho, Kang-Jin; Hwang, Young-Soo; Park, Ro-Dong

CS Department of Biochemistry, National Institute of Agricultural Science and Technology, S. Korea

SO Han'guk Nonghwa Hakhoechi (1999), 42(3), 262-266

CODEN: JKACA7; ISSN: 0368-2897

PB Korean Society of Agricultural Chemistry and Biotechnology

DT Journal

LA Korean

L3 ANSWER 3 OF 3 CA COPYRIGHT 2005 ACS on STN

AB Diterpenoids from roots of **Agastache rugosa** were studied. The dried roots were extracted with 95% EtOH, the EtOH extract was partitioned between **chloroform** and water, the **chloroform** layer was subjected to silica gel CC, using **n-hexane** with gradually increasing proportion of EtOAc as eluent, after further separation agastol (I), a new diterpene, with its isomer named isoagastol (II) were obtained. Their structures were established on the basis of spectral methods and identified as 11,14-dihydroxy-12-methoxy-19(4 \rightarrow 3)abeo-abieta-4(18),8,11,13-tetraen-7-one for agastol and 11,14-dihydroxy-12-methoxy-19(4 \rightarrow 3)abeo-abieta-3,8,11,13-tetraen-7-one, for isoagastol. Isoagastol was isolated for the first time from natural sources.

AN 130:122169 CA

TI Diterpenoids from roots of **Agastache rugosa**

AU Zou, Zhongmei; Yu, Deququan; Cong, Puzhu

CS Institute of Materia Medica, Chinese Academy of Medical Sciences, Beijing, 100050, Peop. Rep. China

SO Journal of Chinese Pharmaceutical Sciences (1997), 6(3), 115-118

CODEN: JCHSE4; ISSN: 1003-1057

PB Beijing Medical University, School of Pharmaceutical Sciences

DT Journal

LA English

RE.CNT 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=>

=> s indigofera tinctoria

L4 127 INDIGOFERA TINCTORIA

=> s dehydrocholic acid
L5 1516 DEHYDROCHOLIC ACID

=> s l4 and l5
L6 1 L4 AND L5

=> d

L6 ANSWER 1 OF 1 CA COPYRIGHT 2005 ACS on STN
AN 139:240385 CA
TI Pharmaceutical composition and process for isolation of
trans-tetracos-15-enoic acid and method of treatment for hepatotoxicity
IN Handa, Sukhdev Swami; Singh, Bupinder; Chandan, Bal Krishan; Saxena, Ajit
Kumar; Bhardwaj, Vikram; Gupta, V. N.; Suri, Om Parkash; Satti, Naresh
Kumar; Suri, Krishan Avtar
PA India
SO U.S. Pat. Appl. Publ., 15 pp.
CODEN: USXXCO
DT Patent
LA English
FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2003175363	A1	20030918	US 2002-73548	20020211
	US 2003157198	A1	20030821	US 2002-102147	20020320
	WO 2003080081	A2	20031002	WO 2002-IB1212	20020325
	WO 2003080081	A3	20041223		
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,				
	CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,				
	GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,				
	LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH,				
	PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ,				
	UA, UG, US, UZ, VN, YU, ZA, ZM, ZW				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY,				
	KG, KZ, MD, RU, TJ, TM, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB,				
	GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA,				
	GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP	1509285	A2	20050302	EP 2002-722566	20020325
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,				
	IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
PRAI	US 2002-73548	A2	20020211		
	WO 2002-IB1212	W	20020325		

=>
=> s indigo
L7 8179 INDIGO

=> d his

(FILE 'HOME' ENTERED AT 15:46:04 ON 15 AUG 2005)

FILE 'CA, BIOSIS, MEDLINE' ENTERED AT 15:46:17 ON 15 AUG 2005

L1 161 S AGASTACHE RUGOSA OR PATCHOLI
L2 108333 S N-HEXANE OR CHLOROFORM OR N-BUTANOL
L3 3 S L1 AND L2
L4 127 S INDIGOFERA TINCTORIA
L5 1516 S DEHYDROCHOLIC ACID
L6 1 S L4 AND L5
L7 8179 S INDIGO

=> s l7 and l5
L8 1 L7 AND L5

=> d

L8 ANSWER 1 OF 1 CA COPYRIGHT 2005 ACS on STN

AN 65:66754 CA
OREF 65:12468f-g
TI The variable catalytic behavior of various bile acids
AU Krause, A.; Domka, F.; Marciniec, B.
CS Univ. Poznan, Pol.
SO Oesterreichische Chemiker-Zeitung (1966), 67(8), 283-4
CODEN: OCHZA8; ISSN: 0369-7061
DT Journal
LA German

=> d his

(FILE 'HOME' ENTERED AT 15:46:04 ON 15 AUG 2005)

FILE 'CA, BIOSIS, MEDLINE' ENTERED AT 15:46:17 ON 15 AUG 2005

L1 161 S AGASTACHE RUGOSA OR PATCHOLI
L2 108333 S N-HEXANE OR CHLOROFORM OR N-BUTANOL
L3 3 S L1 AND L2
L4 127 S INDIGOFERA TINCTORIA
L5 1516 S DEHYDROCHOLIC ACID
L6 1 S L4 AND L5
L7 8179 S INDIGO
L8 1 S L7 AND L5

=> d l8 ab,bib

L8 ANSWER 1 OF 1 CA COPYRIGHT 2005 ACS on STN
AB The catalytic behavior of cholic acid (I), deoxycholic acid (II), and **dehydrocholic acid** (III) was studied by means of the decolorizing-time of **indigo** carmine. The order of catalytic activity was: I, III, II. An explanation is given concerning the rather good catalytic properties of III in spite of the absence of OH groups which are responsible for the activity of I.

AN 65:66754 CA
OREF 65:12468f-g
TI The variable catalytic behavior of various bile acids
AU Krause, A.; Domka, F.; Marciniec, B.
CS Univ. Poznan, Pol.
SO Oesterreichische Chemiker-Zeitung (1966), 67(8), 283-4
CODEN: OCHZA8; ISSN: 0369-7061
DT Journal
LA German

=> file reg

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STRUCTURE FILE UPDATES: 14 AUG 2005 HIGHEST RN 860111-75-7
DICTIONARY FILE UPDATES: 14 AUG 2005 HIGHEST RN 860111-75-7

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*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added,   *
* effective March 20, 2005. A new display format, IDERL, is now    *
* available and contains the CA role and document type information. *
*
*****
```

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:

<http://www.cas.org/ONLINE/DBSS/registryss.html>

=> s indigo/cn

L9 2 INDIGO/CN

=> d 1-2

L9 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2005 ACS on STN

RN 68651-46-7 REGISTRY

ED Entered STN: 16 Nov 1984

CN Indigo (dye) (9CI) (CA INDEX NAME)

OTHER NAMES:

CN Ai

CN **Indigo**

CN Nile

MF Unspecified

CI COM, MAN

LC STN Files: AGRICOLA, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CEN, CIN, EMBASE, PIRA, PROMT, TOXCENTER, USPAT2, USPATFULL

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

553 REFERENCES IN FILE CA (1907 TO DATE)

16 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

554 REFERENCES IN FILE CAPLUS (1907 TO DATE)

L9 ANSWER 2 OF 2 REGISTRY COPYRIGHT 2005 ACS on STN

RN 482-89-3 REGISTRY

ED Entered STN: 16 Nov 1984

CN 3H-Indol-3-one, 2-(1,3-dihydro-3-oxo-2H-indol-2-ylidene)-1,2-dihydro-(9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN Indigo Pure BASF (6CI)

OTHER NAMES:

CN Δ2,2'-Bipseudoindoxyl

CN 11669 Blue

CN Blue No. 201

CN C Blue 22

CN C.I. 73000

CN C.I. Natural Blue 1

CN C.I. Pigment Blue 66

CN C.I. Vat Blue 1

CN Cystoceva

CN D and C Blue No. 6

CN D&C Blue No. 6

CN D+C Blue No. 6

CN Diindogen

CN **Indigo**

CN Indigo Blue

CN Indigo Ciba

CN Indigo Ciba SL
 CN Indigo J
 CN Indigo N
 CN Indigo NAC
 CN Indigo NACCO
 CN Indigo P
 CN Indigo PLN
 CN Indigo Powder W
 CN Indigo Pure BASF Powder K
 CN Indigo Synthetic
 CN Indigo VS
 CN Indigotin
 CN Indigotin (natural)
 CN Indigotine
 CN Japan Blue 201
 CN Lithosol Deep Blue B
 CN Mitsui Indigo Paste
 CN Mitsui Indigo Pure
 CN Mitsui Indigo Pure EXN
 CN Monolite Fast Navy Blue BV
 CN Natural Blue 1
 CN Natural blue indigotin
 CN Pigment Blue 66
 CN Pigment Indigo
 CN Pigment Indigo V
 CN Reduced Dark Blue VB
 CN Synthetic Indigo
 CN Synthetic Indigo TS
 CN Vat Blue 1
 CN Vat Dark Blue VB
 CN Vulcafix Blue R
 CN Vulcafor Blue A
 CN Vulcanosine Dark Blue L

ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for DISPLAY

DR 12000-74-7, 12626-73-2, 11129-41-2, 93660-98-1, 136797-30-3, 210488-46-3

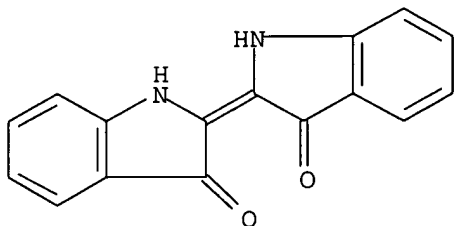
MF C16 H10 N2 O2

CI COM

LC STN Files: AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOBUSINESS, BIOSIS,
 BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS,
 CHEMLIST, CIN, CSCHM, CSNB, DDFU, DETHERM*, DRUGU, EMBASE, GMELIN*,
 HODOC*, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS,
 NIOSHTIC, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, USPAT2, USPATFULL
 (*File contains numerically searchable property data)

Other Sources: DSL**, EINECS**, TSCA**

(**Enter CHEMLIST File for up-to-date regulatory information)



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1424 REFERENCES IN FILE CA (1907 TO DATE)

48 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

1426 REFERENCES IN FILE CAPLUS (1907 TO DATE)

18 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> s indigo carmine and indigofera tinctoria

536 INDIGO
94 CARMINE
7 INDIGO CARMINE
 (INDIGO (W) CARMINE)
15 INDIGOFERA
33 TINCTORIA
1 INDIGOFERA TINCTORIA
 (INDIGOFERA (W) TINCTORIA)
L10 0 INDIGO CARMINE AND INDIGOFERA TINCTORIA